



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:38 PM UTC

PDB ID : 4ZV9 / pdb_00004zv9
Title : 2.00 Angstrom resolution crystal structure of an uncharacterized protein from Escherichia coli O157:H7 str. Sakai
Authors : Halavaty, A.S.; Wawrzak, Z.; Filippova, E.V.; Kiryukhina, O.; Endres, M.; Joachimiak, A.; Anderson, W.F.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2015-05-18
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

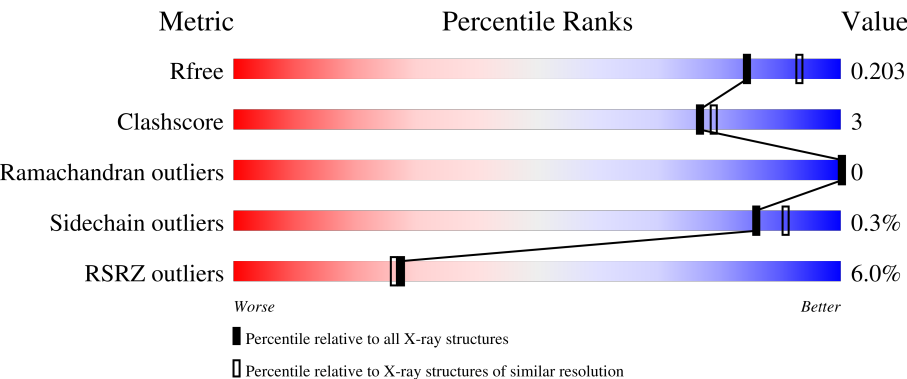
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



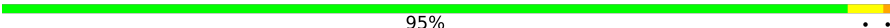
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	238	2% 97% ..
1	B	238	95% 5%
1	C	238	31% 94% 5% .
1	D	238	2% 94% 6%

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Mol	Chain	Length	Quality of chain
1	E	238	 95% . .
1	F	238	 94% 6%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12803 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	S	Se	0	6	0
			1910	1218	326	358	3	5			
1	B	238	Total	C	N	O	S	Se	0	7	0
			1926	1228	336	355	3	4			
1	C	238	Total	C	N	O	S	Se	0	0	0
			1856	1187	315	348	3	3			
1	D	238	Total	C	N	O	S	Se	0	5	0
			1898	1211	321	358	3	5			
1	E	238	Total	C	N	O	S	Se	0	5	0
			1900	1212	322	356	5	5			
1	F	238	Total	C	N	O	S	Se	0	4	0
			1892	1207	325	353	3	4			

There are 18 discrepancies between the modelled and reference sequences:

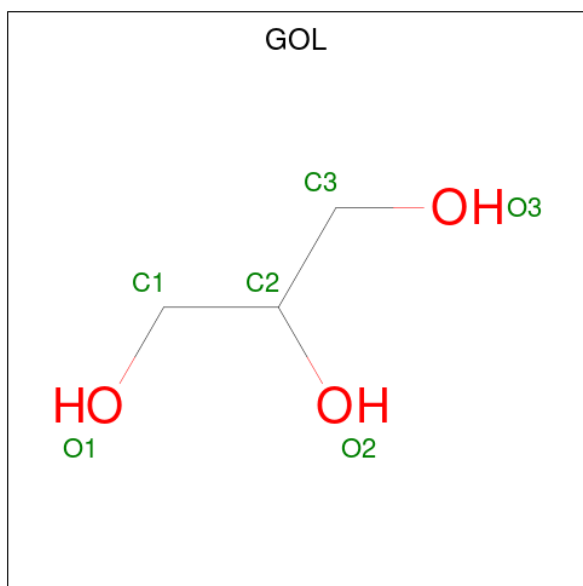
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q7AAT4
A	-1	ASN	-	expression tag	UNP Q7AAT4
A	0	ALA	-	expression tag	UNP Q7AAT4
B	57	SER	-	expression tag	UNP Q7AAT4
B	58	ASN	-	expression tag	UNP Q7AAT4
B	59	ALA	-	expression tag	UNP Q7AAT4
C	57	SER	-	expression tag	UNP Q7AAT4
C	58	ASN	-	expression tag	UNP Q7AAT4
C	59	ALA	-	expression tag	UNP Q7AAT4
D	57	SER	-	expression tag	UNP Q7AAT4
D	58	ASN	-	expression tag	UNP Q7AAT4
D	59	ALA	-	expression tag	UNP Q7AAT4
E	57	SER	-	expression tag	UNP Q7AAT4
E	58	ASN	-	expression tag	UNP Q7AAT4
E	59	ALA	-	expression tag	UNP Q7AAT4
F	57	SER	-	expression tag	UNP Q7AAT4
F	58	ASN	-	expression tag	UNP Q7AAT4

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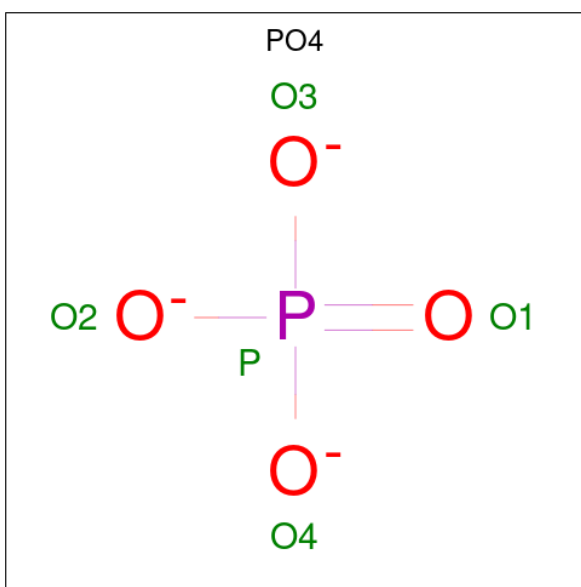
Chain	Residue	Modelled	Actual	Comment	Reference
F	59	ALA	-	expression tag	UNP Q7AAT4

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



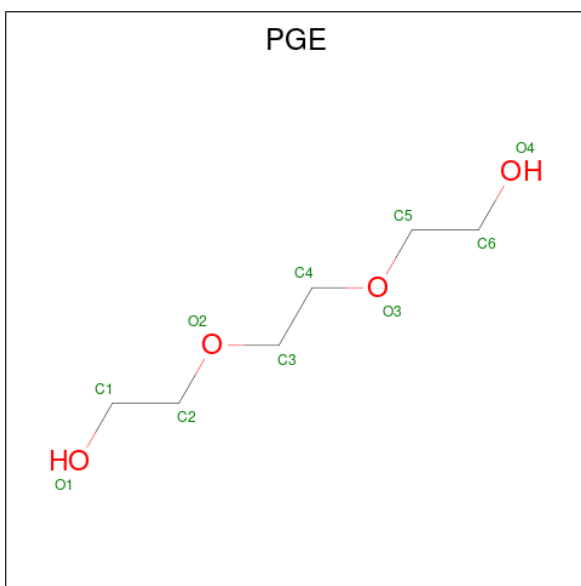
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



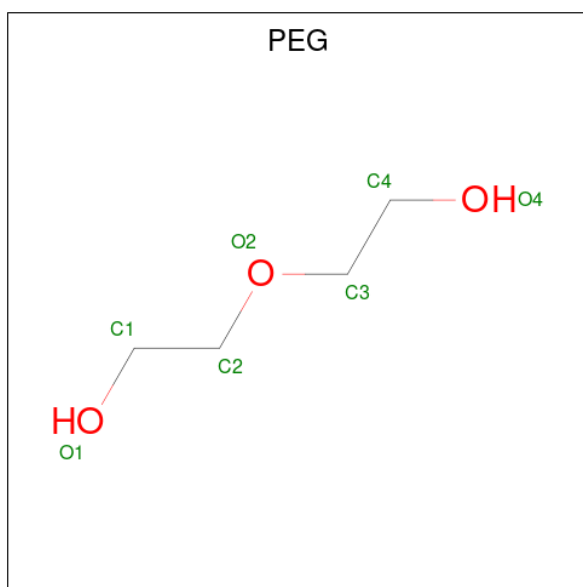
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	C	O	0	0
			7	4	3		

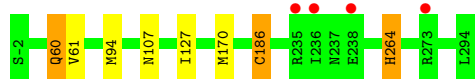
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	183	Total	O	0	4
			184	184		
6	B	333	Total	O	0	16
			341	341		
6	C	50	Total	O	0	0
			50	50		
6	D	142	Total	O	0	5
			147	147		
6	E	295	Total	O	0	6
			299	299		
6	F	309	Total	O	0	12
			318	318		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

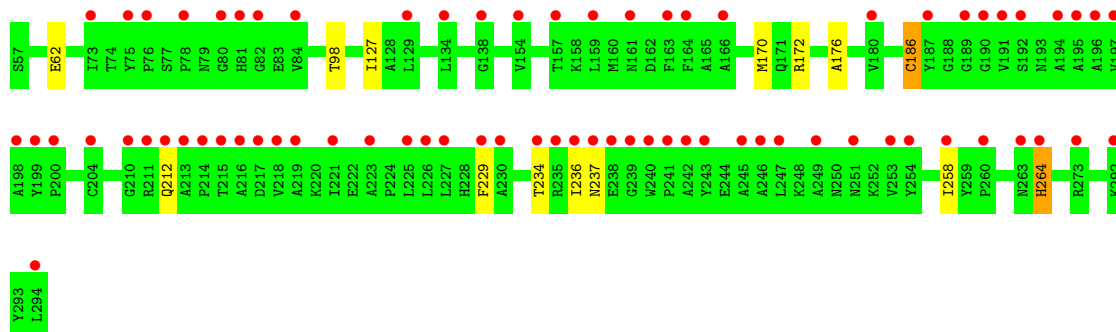
- Molecule 1: Uncharacterized protein



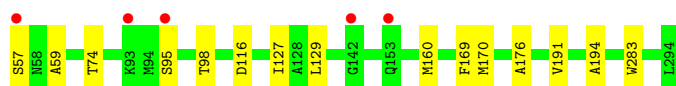
- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



- Molecule 1: Uncharacterized protein



● Molecule 1: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.61Å 96.34Å 129.62Å 90.00° 98.22° 90.00°	Depositor
Resolution (Å)	38.79 – 2.00 38.79 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.0 (38.79-2.00) 92.0 (38.79-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.168 , 0.199 (Not available) , 0.203	Depositor DCC
R_{free} test set	5263 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12803	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, PGE, CME, GOL, PEG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.57	0/1949	0.77	1/2644 (0.0%)
1	B	0.63	0/1965	0.78	0/2662
1	C	0.53	0/1895	0.80	1/2574 (0.0%)
1	D	0.53	0/1937	0.77	0/2629
1	E	0.65	0/1929	0.80	1/2618 (0.0%)
1	F	0.62	0/1931	0.77	0/2620
All	All	0.59	0/11606	0.78	3/15747 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	58	ASN	N-CA-C	6.24	116.81	108.38
1	C	264	HIS	N-CA-CB	-5.63	100.73	109.69
1	A	264	HIS	N-CA-CB	-5.42	101.07	109.69

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1910	0	1834	7	0
1	B	1926	0	1863	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1856	0	1783	8	0
1	D	1898	0	1814	14	0
1	E	1900	0	1820	14	0
1	F	1892	0	1819	13	0
2	A	18	0	24	0	0
2	B	6	0	8	0	0
2	D	6	0	8	0	0
2	E	12	0	16	0	0
2	F	18	0	24	0	0
3	B	5	0	0	0	0
4	E	10	0	14	3	0
5	F	7	0	10	0	0
6	A	184	0	0	2	0
6	B	341	0	0	8	0
6	C	50	0	0	0	0
6	D	147	0	0	2	0
6	E	299	0	0	3	0
6	F	318	0	0	7	0
All	All	12803	0	11037	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (65) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:129:LEU:HD22	1:D:170[B]:MSE:HE3	1.42	1.02
1:D:169:PHE:CD2	1:D:170[B]:MSE:HE2	2.01	0.94
1:E:94[A]:MSE:HE2	6:E:508:HOH:O	1.76	0.82
1:D:169:PHE:HD2	1:D:170[B]:MSE:HE2	1.47	0.79
1:D:95:SER:HB3	6:D:496:HOH:O	1.87	0.73
1:D:169:PHE:CE2	1:D:170[B]:MSE:HE2	2.23	0.73
1:C:212:GLN:NE2	1:C:236:ILE:O	2.21	0.73
1:A:94:MSE:HE2	6:A:457:HOH:O	1.88	0.71
1:A:186:CME:SG	1:A:264:HIS:NE2	2.59	0.71
1:F:212:GLN:HG2	6:F:568:HOH:O	1.90	0.70
4:E:303:PGE:H2	1:F:174:PRO:HD3	1.75	0.69
1:B:62:GLU:OE2	1:E:57:SER:HA	1.93	0.68
4:E:303:PGE:C2	1:F:174:PRO:HD3	2.22	0.68
1:D:129:LEU:HD22	1:D:170[B]:MSE:CE	2.22	0.65
1:B:60:GLN:OE1	1:E:58:ASN:HB3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:160[B]:MSE:HE3	6:E:632:HOH:O	1.97	0.65
1:A:60[A]:GLN:HG3	1:A:61:VAL:N	2.11	0.64
1:C:229:PHE:HB2	1:C:258:ILE:HD13	1.81	0.62
1:C:212:GLN:OE1	1:C:212:GLN:N	2.31	0.62
1:F:171:GLN:NE2	6:F:402:HOH:O	2.33	0.60
1:B:211[A]:ARG:NH2	6:B:404:HOH:O	2.34	0.60
1:F:94:MSE:HE2	6:F:466:HOH:O	2.01	0.60
1:B:94:MSE:HE2	6:B:508:HOH:O	2.02	0.59
1:C:186:CME:SG	1:C:264:HIS:NE2	2.61	0.59
1:B:186:CME:HB3	6:B:622[B]:HOH:O	2.02	0.58
1:F:160[B]:MSE:HE3	6:F:675:HOH:O	2.04	0.56
1:E:160[B]:MSE:HE2	1:E:191:VAL:HG22	1.86	0.56
1:F:108:ARG:NH1	6:F:404:HOH:O	2.39	0.55
6:B:643:HOH:O	1:E:57:SER:HB2	2.07	0.55
1:B:288[A]:LYS:HE2	6:B:545:HOH:O	2.07	0.54
4:E:303:PGE:H22	1:F:174:PRO:HD3	1.90	0.54
1:B:60:GLN:OE1	1:E:58:ASN:CB	2.55	0.54
1:D:160[B]:MSE:HE3	1:D:191:VAL:HG22	1.91	0.53
1:A:107:ASN:HB3	1:A:186:CME:HZ2	1.90	0.53
1:C:234:THR:HA	1:C:237:ASN:OD1	2.08	0.53
1:D:283:TRP:CG	6:D:468:HOH:O	2.63	0.52
1:F:97:LYS:NZ	6:F:405:HOH:O	2.41	0.52
1:B:288[A]:LYS:HD2	6:B:660:HOH:O	2.09	0.51
1:D:160[B]:MSE:HE2	1:D:194:ALA:HB3	1.95	0.49
1:E:91:PRO:HG2	1:E:94[B]:MSE:HE2	1.94	0.49
1:A:60[A]:GLN:NE2	6:A:401[A]:HOH:O	2.32	0.48
1:B:148[B]:ARG:HD2	6:B:650:HOH:O	2.16	0.45
1:E:186[B]:CME:HB2	1:E:186[B]:CME:HE3	1.47	0.45
1:B:186:CME:HZ2	1:B:265:GLY:HA2	1.98	0.45
1:A:127:ILE:HG21	1:A:170[B]:MSE:HE1	1.99	0.45
1:E:127:ILE:HG21	1:E:170:MSE:HE1	1.99	0.45
1:A:170[A]:MSE:HE2	1:A:170[A]:MSE:HA	1.99	0.44
1:D:127:ILE:HG21	1:D:170[A]:MSE:HE1	1.99	0.44
1:B:186:CME:HZ2	1:B:265:GLY:CA	2.48	0.43
1:B:127:ILE:HG21	1:B:170:MSE:HE1	2.00	0.43
1:E:91:PRO:HB2	1:E:94[B]:MSE:HG2	2.01	0.43
6:B:561:HOH:O	1:E:58:ASN:HB2	2.18	0.43
1:F:74:THR:HG23	6:F:631:HOH:O	2.19	0.43
1:E:186[A]:CME:HB2	6:E:549[A]:HOH:O	2.18	0.42
1:F:127:ILE:HG21	1:F:170:MSE:HE1	2.02	0.42
1:B:98:THR:O	1:B:176:ALA:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:ILE:HG21	1:C:170:MSE:HE1	2.00	0.42
1:F:98:THR:O	1:F:176:ALA:HA	2.19	0.42
1:D:169:PHE:CE2	1:D:170[B]:MSE:CE	3.00	0.41
1:C:62:GLU:OE1	1:D:57:SER:HA	2.21	0.41
1:D:98:THR:O	1:D:176:ALA:HA	2.21	0.41
1:C:98:THR:O	1:C:176:ALA:HA	2.20	0.41
1:F:186:CME:HZ2	1:F:273[B]:ARG:HE	1.85	0.41
1:D:59:ALA:HB1	1:D:116:ASP:OD1	2.21	0.41
1:E:98:THR:O	1:E:176:ALA:HA	2.21	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	241/238 (101%)	236 (98%)	5 (2%)	0	100	100
1	B	242/238 (102%)	237 (98%)	5 (2%)	0	100	100
1	C	235/238 (99%)	230 (98%)	5 (2%)	0	100	100
1	D	240/238 (101%)	234 (98%)	6 (2%)	0	100	100
1	E	239/238 (100%)	234 (98%)	5 (2%)	0	100	100
1	F	239/238 (100%)	232 (97%)	7 (3%)	0	100	100
All	All	1436/1428 (101%)	1403 (98%)	33 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	191/182 (105%)	189 (99%)	2 (1%)	68	75
1	B	192/182 (106%)	192 (100%)	0	100	100
1	C	185/182 (102%)	184 (100%)	1 (0%)	81	87
1	D	190/182 (104%)	189 (100%)	1 (0%)	81	87
1	E	189/182 (104%)	189 (100%)	0	100	100
1	F	189/182 (104%)	189 (100%)	0	100	100
All	All	1136/1092 (104%)	1132 (100%)	4 (0%)	86	89

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60[A]	GLN
1	A	60[B]	GLN
1	C	172	ARG
1	D	74	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	111	ASN
1	D	111	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

7 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CME	F	186	1	8,9,10	0.82	0	6,9,11	1.04	0
1	CME	A	186	1	8,9,10	0.84	0	6,9,11	1.61	1 (16%)
1	CME	D	186	1	8,9,10	0.78	0	6,9,11	0.77	0
1	CME	E	186[B]	1	8,9,10	0.85	0	6,9,11	1.24	1 (16%)
1	CME	B	186	1	8,9,10	1.61	1 (12%)	6,9,11	2.63	3 (50%)
1	CME	E	186[A]	1	8,9,10	0.93	0	6,9,11	0.87	0
1	CME	C	186	1	8,9,10	0.71	0	6,9,11	1.45	1 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CME	F	186	1	-	3/5/8/10	-
1	CME	A	186	1	-	3/5/8/10	-
1	CME	D	186	1	-	2/5/8/10	-
1	CME	E	186[B]	1	-	1/5/8/10	-
1	CME	B	186	1	-	4/5/8/10	-
1	CME	E	186[A]	1	-	3/5/8/10	-
1	CME	C	186	1	-	4/5/8/10	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	186	CME	CB-SG	-4.11	1.68	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	186	CME	CB-SG-SD	-4.38	92.53	103.86
1	B	186	CME	CA-CB-SG	-3.20	101.31	114.45
1	A	186	CME	CB-CA-C	2.92	118.72	110.80
1	E	186[B]	CME	CB-SG-SD	-2.83	96.55	103.86
1	C	186	CME	CB-CA-C	2.81	118.44	110.80
1	B	186	CME	CB-CA-C	-2.65	103.61	110.80

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	186	CME	N-CA-CB-SG
1	B	186	CME	SD-CE-CZ-OH
1	F	186	CME	CZ-CE-SD-SG
1	B	186	CME	CE-SD-SG-CB
1	E	186[A]	CME	CE-SD-SG-CB
1	F	186	CME	CE-SD-SG-CB
1	E	186[B]	CME	CE-SD-SG-CB
1	A	186	CME	SD-CE-CZ-OH
1	C	186	CME	SD-CE-CZ-OH
1	A	186	CME	N-CA-CB-SG
1	C	186	CME	N-CA-CB-SG
1	B	186	CME	CZ-CE-SD-SG
1	C	186	CME	CZ-CE-SD-SG
1	E	186[A]	CME	CZ-CE-SD-SG
1	D	186	CME	CA-CB-SG-SD
1	A	186	CME	CZ-CE-SD-SG
1	C	186	CME	CA-CB-SG-SD
1	D	186	CME	SD-CE-CZ-OH
1	E	186[A]	CME	CA-CB-SG-SD
1	F	186	CME	CA-CB-SG-SD

There are no ring outliers.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	F	186	CME	1	0
1	A	186	CME	2	0
1	E	186[B]	CME	1	0
1	B	186	CME	3	0
1	E	186[A]	CME	1	0
1	C	186	CME	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

13 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	F	303	-	5,5,5	0.18	0	5,5,5	0.39	0
2	GOL	E	302	-	5,5,5	0.30	0	5,5,5	0.92	0
2	GOL	F	302	-	5,5,5	0.43	0	5,5,5	1.02	0
4	PGE	E	303	-	9,9,9	0.62	0	8,8,8	0.71	0
2	GOL	A	303	-	5,5,5	0.29	0	5,5,5	0.40	0
2	GOL	A	301	-	5,5,5	0.37	0	5,5,5	0.36	0
2	GOL	B	302	-	5,5,5	0.11	0	5,5,5	0.60	0
2	GOL	E	301	-	5,5,5	0.31	0	5,5,5	0.41	0
3	PO4	B	301	-	4,4,4	1.02	0	6,6,6	0.40	0
2	GOL	F	301	-	5,5,5	0.22	0	5,5,5	0.38	0
2	GOL	A	302	-	5,5,5	0.32	0	5,5,5	0.26	0
5	PEG	F	304	-	6,6,6	0.50	0	5,5,5	0.25	0
2	GOL	D	301	-	5,5,5	0.29	0	5,5,5	0.51	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	F	303	-	-	2/4/4/4	-
2	GOL	E	302	-	-	4/4/4/4	-
2	GOL	F	302	-	-	0/4/4/4	-
4	PGE	E	303	-	-	3/7/7/7	-
2	GOL	A	303	-	-	2/4/4/4	-
2	GOL	A	301	-	-	2/4/4/4	-
2	GOL	B	302	-	-	4/4/4/4	-
2	GOL	E	301	-	-	1/4/4/4	-
2	GOL	F	301	-	-	3/4/4/4	-
2	GOL	A	302	-	-	0/4/4/4	-
5	PEG	F	304	-	-	2/4/4/4	-
2	GOL	D	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	GOL	C1-C2-C3-O3
2	A	303	GOL	O1-C1-C2-C3
2	E	302	GOL	O1-C1-C2-O2
2	F	301	GOL	C1-C2-C3-O3
2	F	301	GOL	O2-C2-C3-O3
2	B	302	GOL	O2-C2-C3-O3
2	E	302	GOL	O2-C2-C3-O3
4	E	303	PGE	O1-C1-C2-O2
2	B	302	GOL	C1-C2-C3-O3
2	E	302	GOL	O1-C1-C2-C3
2	E	302	GOL	C1-C2-C3-O3
2	F	303	GOL	C1-C2-C3-O3
2	A	301	GOL	O2-C2-C3-O3
5	F	304	PEG	O2-C3-C4-O4
5	F	304	PEG	O1-C1-C2-O2
2	F	303	GOL	O2-C2-C3-O3
2	A	303	GOL	O1-C1-C2-O2
2	E	301	GOL	O2-C2-C3-O3
2	F	301	GOL	O1-C1-C2-O2
4	E	303	PGE	O3-C5-C6-O4
2	B	302	GOL	O1-C1-C2-C3
2	B	302	GOL	O1-C1-C2-O2
4	E	303	PGE	O2-C3-C4-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	303	PGE	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	234/238 (98%)	0.17	4 (1%) 69 68	14, 33, 58, 94	4 (1%)
1	B	234/238 (98%)	-0.52	0 100 100	12, 21, 34, 80	6 (2%)
1	C	234/238 (98%)	1.48	73 (31%) 1 1	33, 76, 130, 144	0
1	D	234/238 (98%)	0.26	5 (2%) 63 63	16, 42, 63, 129	3 (1%)
1	E	234/238 (98%)	-0.43	1 (0%) 88 88	13, 23, 38, 79	2 (0%)
1	F	234/238 (98%)	-0.50	1 (0%) 88 88	12, 22, 38, 84	3 (1%)
All	All	1404/1428 (98%)	0.08	84 (5%) 27 26	12, 30, 92, 144	18 (1%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	57	SER	4.8
1	C	237	ASN	4.3
1	C	214	PRO	4.1
1	E	57	SER	4.0
1	C	219	ALA	3.7
1	C	247	LEU	3.7
1	C	78	PRO	3.7
1	C	80	GLY	3.6
1	C	239	GLY	3.6
1	C	235	ARG	3.6
1	C	236	ILE	3.6
1	C	200	PRO	3.5
1	A	235	ARG	3.5
1	C	196	ALA	3.4
1	C	213	ALA	3.4
1	C	245	ALA	3.3
1	C	243	TYR	3.2
1	C	240	TRP	3.2
1	C	249	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	95	SER	3.1
1	C	241	PRO	3.1
1	C	129	LEU	3.1
1	A	236	ILE	3.1
1	C	84	VAL	3.0
1	C	159	LEU	3.0
1	C	192	SER	2.9
1	C	242	ALA	2.9
1	C	190	GLY	2.9
1	C	225	LEU	2.8
1	A	273[A]	ARG	2.8
1	C	198	ALA	2.8
1	C	254	TYR	2.8
1	C	154	VAL	2.8
1	C	197	VAL	2.8
1	C	221	ILE	2.8
1	C	166	ALA	2.7
1	C	216	ALA	2.7
1	C	75	TYR	2.7
1	C	238	GLU	2.7
1	C	76	PRO	2.7
1	C	263	ASN	2.6
1	D	93	LYS	2.6
1	C	223	ALA	2.6
1	C	163	PHE	2.6
1	C	264	HIS	2.6
1	C	251	ASN	2.6
1	C	211	ARG	2.6
1	D	153	GLN	2.6
1	C	189	GLY	2.6
1	C	273	ARG	2.5
1	C	230	ALA	2.5
1	C	180	VAL	2.5
1	C	191	VAL	2.5
1	C	218	VAL	2.4
1	C	229	PHE	2.4
1	C	253	VAL	2.4
1	C	258	ILE	2.4
1	A	238	GLU	2.4
1	C	164	PHE	2.4
1	C	157	THR	2.4
1	C	227	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	226	LEU	2.3
1	C	217	ASP	2.3
1	C	81	HIS	2.3
1	C	138	GLY	2.3
1	C	210	GLY	2.3
1	C	194	ALA	2.3
1	C	246	ALA	2.3
1	C	199	TYR	2.3
1	C	161	ASN	2.2
1	C	204	CYS	2.2
1	C	234	THR	2.2
1	D	142	GLY	2.2
1	C	215	THR	2.2
1	C	294	LEU	2.1
1	C	187	TYR	2.1
1	C	82	GLY	2.1
1	C	292	LYS	2.1
1	C	195	ALA	2.1
1	C	73	ILE	2.1
1	C	134	LEU	2.0
1	C	212	GLN	2.0
1	F	93	LYS	2.0
1	C	260	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	CME	C	186	10/11	0.69	0.20	71,78,91,92	0
1	CME	E	186[A]	10/11	0.73	0.22	19,32,46,52	10
1	CME	E	186[B]	10/11	0.73	0.22	18,27,34,34	10
1	CME	A	186	10/11	0.88	0.15	29,42,71,79	0
1	CME	B	186	10/11	0.89	0.14	16,33,46,57	0
1	CME	D	186	10/11	0.94	0.10	27,38,66,77	0
1	CME	F	186	10/11	0.94	0.13	16,25,65,75	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GOL	A	301	6/6	0.73	0.15	73,75,78,79	0
2	GOL	E	301	6/6	0.75	0.16	62,73,77,86	0
3	PO4	B	301	5/5	0.75	0.15	48,49,54,63	5
4	PGE	E	303	10/10	0.75	0.20	48,60,63,65	0
2	GOL	D	301	6/6	0.80	0.15	45,51,57,63	0
5	PEG	F	304	7/7	0.81	0.16	52,55,62,63	0
2	GOL	F	301	6/6	0.82	0.18	58,63,66,67	0
2	GOL	A	303	6/6	0.83	0.15	51,52,56,57	0
2	GOL	B	302	6/6	0.84	0.18	39,49,59,64	0
2	GOL	F	303	6/6	0.85	0.18	40,47,66,72	0
2	GOL	E	302	6/6	0.85	0.13	37,43,48,50	0
2	GOL	A	302	6/6	0.90	0.10	45,49,50,53	0
2	GOL	F	302	6/6	0.92	0.12	28,35,37,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.